# Solving Linear Programs with Complementarity Constraints using Branch-and-Cut

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Abstract A linear program with linear complementarity constraints (LPCC) requires the minimization of a linear objective over a set of linear constraints together with additional linear complementarity constraints. This class has emerged as a modeling paradigm for a broad collection of problems, including bilevel programs, Stackelberg games, inverse quadratic programs, and problems involving equilibrium constraints. The presence of the complementarity constraints results in a nonconvex optimization problem. We develop a branch-and-cut algorithm to find a global optimum for this class of optimization problems, where we branch directly on complementarities. We develop branching rules and feasibility recovery procedures and demonstrate their computational effectiveness in a comparison with CPLEX. The implementation builds on CPLEX through the use of callback routines. The computational results show that our approach is a strong alternative to constructing an integer programming formulation using big-M terms to represent bounds for variables, with testing conducted on general LPCCs as well as on instances generated from bilevel programs with convex quadratic lower level problems.

Keywords linear programs with complementarity constraints  $\cdot$  MPECs  $\cdot$  branch-and-cut

## 1 Introduction

A linear program with linear complementarity constraints (LPCC), which minimizes a linear objective function over a set of linear constraints with additional linear complementarity constraints, is a nonconvex, disjunctive optimization problem. In §1.1, we present the mathematical formulation of the general LPCC we use throughout this paper. In §1.2, various existing algorithms designed for solving LPCCs are reviewed. Most of these existing methods are only able to obtain a stationary solution and incapable of ascertaining the quality of the solution. This is the major drawback for the existing solvers. In this paper, we mainly focus on finding the global resolution of the LPCC, and we achieve this goal through two steps:

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- Step 1 Study various valid constraints by exploiting the complementarity constraints directly, and evaluate the benefit of these constraints on the value of the linear relaxation of the LPCC. We have previously discussed valid constraints for the LPCC in [49], and we briefly recap these constraints in §2.2.
- Step 2 Propose a branch-and-cut algorithm to globally solve the LPCC problem, where cuts are derived from various valid constraints studied in Step 1 and branching is imposed on the complementarity constraints. A general LPCC solver has been developed based on this branch-and-cut approach, and it is able to compete with the existing MIP-based solvers like CPLEX.

The branch-and-cut algorithm is introduced in  $\S1.3$ , where we also outline the rest of the paper.

## 1.1 Statement of the Problem

We consider a general formulation of the LPCC in the form suggested by Pang and Fukushima [52]. Given vectors and matrices:  $c \in \mathbb{R}^n$ ,  $d \in \mathbb{R}^m$ ,  $b \in \mathbb{R}^k$ ,  $q \in \mathbb{R}^m$ ,  $A \in \mathbb{R}^{k \times n}$ ,  $B \in \mathbb{R}^{k \times m}$ ,  $N \in \mathbb{R}^{m \times n}$  and  $M \in \mathbb{R}^{m \times m}$ , the LPCC is to find  $(x, y, w) \in \mathbb{R}^n \times \mathbb{R}^m \times \mathbb{R}^m$  in order to solve to global optimality

$$\begin{array}{ll} \underset{(x,y,w)}{\text{minimize}} & c^T x + d^T y \\ \text{subject to } Ax + By \ge b \\ & x \ge 0 \\ \text{and} & 0 \le y \perp w := q + Nx + My \ge 0 \end{array}$$
(1)

where  $a \perp b$  denotes perpendicularity between vectors a and b, i.e.,  $a^T b = 0$ . Without the orthogonality condition  $y \perp w$ , the LPCC is a linear program (LP). The global resolution of the LPCC means the generation of a certificate showing that the problem is in one of its 3 possible states: (a) it is infeasible, (b) it is feasible but unbounded below, or (c) it attains a finite optimal solution. Note that problem (1) is equivalent to  $2^m$  linear programs obtained by making each possible assignment for the complementary variables: either  $y_i = 0$  or  $w_i = 0$  for each  $i = 1, \ldots, m$ ; hence, it is not possible for an LPCC to have a finite optimal value that is not attained.

If the feasible regions for y and w are bounded then there exist diagonal matrices  $\Theta^y$  and  $\Theta^w$  with diagonal entries  $\theta^y_i$  and  $\theta^w_i$  and problem (1) can be formulated as a mixed integer program:

$$\begin{array}{ll} \underset{(x,y,z)}{\operatorname{minimize}} & c^T x + d^T y \\ \text{subject to } Ax + By \ge b \\ & x \ge 0 \\ & 0 \le y \le \Theta^y z \\ & 0 \le q + Nx + My \le \Theta^w (\mathbf{1} - z) \\ \text{and} & z \in \{0, 1\}^m \end{array}$$

$$(2)$$

The obvious drawback of this formulation is that in order to find  $\theta_i^y$  and  $\theta_i^w$  we need to compute valid upper bounds of  $y_i$  and  $w_i$ , not to mention such upper bounds may not exist if the feasible regions for y and/or w are unbounded. To avoid this drawback, in this paper, we present a branch-and-cut algorithm which branches on the complementarity constraint directly. Previous work on branching on complementarity constraints includes [11,19,30].

Problem (1) generalizes the standard linear complementarity problem (LCP) [14]:  $0 \le y \perp q + My \ge 0$ , so the LPCC is NP-Hard. Moreover, affine variational constraints also lead to the problem (1) [46]. Applications of the LPCC are surveyed in [33]. Among these applications, complementarity constraints play three principal roles during the modelling process:

- Modelling KKT optimality conditions that must be satisfied by some of the variables. Such applications include hierarchical optimization such as Stackelberg games [55], inverse convex quadratic programs, indefinite quadratic programs [27,31], and cross-validated support vector regression [42, 43].
- 2. Modelling equilibrium constraints. See for example the texts [16,46], the survey article [51], or a recent paper on market equilibrium in electric power markets [24].
- 3. Modelling certain logical conditions that are required by some practical optimization problems. Such applications include non-convex piecewise linear optimization, quantile minimization [53], and  $\ell_0$ -minimization [13,21].

#### 1.2 Previous Work on Solving LPCCs

Research on algorithms for solving an LPCC can be divided into two main areas: one concerns the development of globally convergent algorithms with a guarantee of finding a suitable stationary point; the other concerns the development of exact algorithms for global resolution of an LPCC. See the survey [39] for a more detailed review.

It is noted that the methods which are able to solve a general LCP can also be extended to solve an LPCC by using the so called sequential LCP Method. Such a procedure can be found in detail in [40]. A complementary pivoting algorithm for an LPCC is an extension of a pivoting algorithm for LCP which handles linear complementarity constraints just as the classic simplex algorithm for linear programs. Such algorithms usually perform in this way: start from a feasible solution, maintain feasibility for all iterations and try to improve the objective function. Under certain constraint qualifications, these methods guarantee convergence to a certain stationary solution. The references [17,26,36] study and implement this type of method to solve the general LPCC. Another way to get a stationary point is through a so called regularization framework [54]: construct a sequence of relaxed problems controlled by some parameter, then obtain a sequence of solutions which converge to a stationary point when the parameter goes to the limit. Each regularized relaxed problem is solved by an NLP based algorithm such as an interior point method. One method of regularization is to introduce a positive parameter  $\phi$  and relax the complementarity constraints in problem (1) using either  $\{y, w \ge 0, y^T w \le \phi\}$  or some other approach [13,21]. An alternative is to put a penalty for violation of the complementarity constraints into the objective, and gradually update the penalty to infinity [44]. A homotopy method has also been proposed [57]. The obvious drawback of these methods is that they are incapable of ascertaining the quality of the computed solution.

The methods for global resolution of an LPCC are mainly based on an enumerative scheme. Several branch-and-bound methods have been proposed for solving an LPCC derived from a bilevel linear program. Bard and Moore [10] proposed a pure branch-and-bound method for solving bilevel linear programs. Hansen et al. [29] enhance this branch-and-bound scheme by exploiting the necessary optimality condition of the inner problem. As opposed to a branch-and-bound method, the references [34,38] study alternative ways to solve an LPCC by using a cutting plane method. Audet et al. [6] proposed a branchand-cut algorithm for solving bilevel linear programs. An RLT method for finding a feasible solution to a problem with both binary and complementarity constraints is proposed in [24]. It follows from the results of [8] that an LPCC can be lifted to an equivalent convex optimization problem so it can in principle be solved globally using a convex optimization algorithm; the drawback to this approach is that the convexity is over the cone of completely positive matrices which is hard to work with computationally.

It is noted that most of the existing methods for global resolution of the LPCC presume the LPCC has a finite optimal value, and this limitation was not resolved until the paper [32]. In that paper, the authors proposed a minimax integer programming formulation of the LPCC, and solve this system using a Benders decomposition method. The method was extended to quadratic programs with complementarity constraints in [7]. A branching scheme for determining boundedness of the optimal value of a linear program with a bilinear objective function was proposed in [5].

The success of the Benders decomposition method [7,32] heavily depends on a so called sparsification process. If the sparsification process is not successful, in the worst case it will be necessary to check every piece of the LPCC. In this paper, we alternatively use a specialized branch-and-cut scheme which is a more systematic enumerative process to get the global resolution of the LPCC, and our algorithm is also able to characterize infeasible and unbounded LPCC problems as well as solve problems with finite optimal value. Moreover we also discuss various valid constraints for the LPCC by exploiting the complementarity structure; this topic has not been fully exploited in the literature for studying the LPCC.

The complementarity structure of an LPCC can be generalized to SOS1 constraints, a type of special ordered set constraint requiring that at most one of a set of variables is nonzero. Recent work on branchand-cut approaches to problems with SOS1 constraints include [20,22]. De Farias et al. [20] considered problems where all the coefficients are nonnegative and their emphasis is on possible families of cutting planes using a sequential lifting procedure. Fischer and Pfetsch [22] emphasize cuts and branching techniques for problems with overlapping SOS1 constraints, that is, sets of complementarity constraints that have variables in common; this structure can be represented with conflict graphs and can be exploited in the derivation of valid cutting planes and in the construction of sophisticated branching rules building on the ideas of Beale and Tomlin [11]. In our formulation, each variable appears in at most one complementarity constraint, so the nice techniques of Fischer and Pfetsch would not be helpful.

#### 1.3 LPCC using Branch-and-Cut

In this paper, we propose a branch-and-cut algorithm for solving the general LPCC problem (1). In §2, we describe the preprocessing phase of our algorithm: in §2.1, a heuristic feasibility recovery procedure is developed to recover a feasible solution of the LPCC which provides a valid upper bound of the LPCC; and in §2.2, the strategy of generating and selecting from various types of cutting planes we studied in [49] is discussed, which could sharpen the LP relaxation and improve the initial lower bound of LPCC. In §3, we present the second phase of our algorithm: branch-and-bound. Various node selection strategies and branching complementarity selection strategies are discussed in §3.1 and §3.2. Our proposed algorithm is able to characterize infeasible and unbounded LPCC problems as well as solve problems with finite optimal value. The algorithm is summarized in §4. In §5, we show the computational results of our branch-and-cut algorithm on solving randomly generated LPCC instances.

In the MIP formulation (2), the binary vector z is only used to model the complementary relationship of the LPCC, and except for the complementarity constraints it does not interact with x and y at all. This observation motivates us to enforce the complementarities through a specialized branching scheme, i.e., branch on complementarities directly without introducing the binary vector z. This kind of specialized branching approach has been studied to solve several problems such as generalized assignment problems [18], nonconvex quadratic programs [56], nonconvex piecewise linear optimization problems [41], and problems with overlapping SOS1 constraints [20,22]. The obvious advantage of using a specialized branching approach for solving the LPCC is that we no longer need  $\theta$  in the formulation, and therefore this approach is also applicable for the case when y or w is unbounded. In fact, even if we know such a  $\theta$ exists, the cost of computing a valid  $\theta$  could be very expensive especially when m is very large. Moreover, introducing the binary vector z will lead to an increase in both the number of variables and the number of constraints, and these Big-M type constraints are usually not tight which will lead to a number of violated complementarities in the solution of the relaxation.

# 2 Preprocessing Phase

When the initial LP relaxation is bounded below, the preprocessing phase will be invoked, consisting of a feasibility recovery process and a cutting plane selection and management process. The feasibility recovery process may provide a valid upper bound for the LPCC, while the cutting plane selection and management process may provide a better lower bound for the LPCC. Both processes may provide a good starting point for the second phase of our algorithm: branch-and-bound.

# 2.1 LPCC Feasibility Recovery

Finding a good feasible solution to an LPCC is an essential component of our branch-and-cut algorithm for globally resolving the LPCC. A good upper bound can help prune nodes quickly, and avoid unnecessary branching. Notice that here we assume the initial LP relaxation is bounded when we apply our feasibility recovery procedures. Our feasibility recovery procedures have some similarities to feasibility pumps for MIP and MINLP [23].

For ease of discussion, we first introduce some notation and definitions.

**Definition 1** Given any binary vector z with dimension m, we define the linear program LPCC(z) as follows:

$$\begin{array}{l} \underset{(x,y)}{\operatorname{minimize}} \quad c^{T} x + d^{T} y \\ \text{subject to } Ax + By \ge b \\ x \ge 0 \\ 0 \le y \\ 0 \le q + Nx + My \\ 0 \ge y_{i} \\ 0 \ge y_{i} \\ 0 \ge (q + Nx + My)_{i} \quad \text{if } z_{i} = 0 \\ 0 \ge (q + Nx + My)_{i} \quad \text{if } z_{i} = 1. \end{array}$$

$$(3)$$

LPCC(z) is a so-called piece of the LPCC corresponding to the binary vector z.

**Definition 2** The feasibility gap of the piece of an LPCC corresponding to the binary vector z, denoted by FG-LPCC(z), is the optimal value of the following linear program:

$$\begin{array}{ll} \underset{(x,y,w)}{\text{minimize}} & (\mathbf{1} - z)^T y + z^T w\\ \text{subject to } Ax + By \ge b\\ & x \ge 0\\ & 0 \le y\\ & 0 \le w := q + Nx + My \end{array}$$

$$(4)$$

where **1** is the vector with all components equal to 1.

Based on the above two definitions, it is obvious that the following proposition is true:

**Proposition 1** LPCC(z) is feasible if and only if FG-LPCC(z)=0

**Definition 3** Binary vectors z and z' are *adjacent* if there is exactly one component that is different between z and z'.

**Definition 4** If binary vectors z and z' are adjacent and FG-LPCC(z) < FG-LPCC(z'), then  $\Delta z = z - z'$  is a *feasibility gap descent direction* for z'.

Just like mixed integer programs, it is often a good idea to recover a feasible solution based on the LP relaxation solution. The most intuitive recovery process is to round the LP relaxation solution into a solution that satisfies all the complementarity constraints. We use this rounding procedure to initialize a new local search feasibility recovery process, detailed in Procedure 1. Notice that we define search *breadth* as the number of candidates that we are going to select from binary vectors which are adjacent to the initial  $z^*$ , and search *depth* as the maximum number of iterations that we are going to perform for each candidate. We can set search *breadth* and search *depth* to control the local search process. The proposed local search procedure can be used to find a feasible solution, although the quality of the recovered feasible solution is not guaranteed. We use optimality based bound tightening [28,47, 58] to resolve this issue, refining the local search feasibility recovery procedure through the addition of the constraints  $lb_{search} \leq c^T x + d^T y \leq ub_{search}$  to (4) when computing the feasibility gap. Procedure 2 describes this refined feasibility recovery process in §5. See Fischer and Pfetsch [22] for primal heuristics that can be used when a variable appears in more than one complementarity constraint.

## 2.2 Cutting Plane Generation and Selection

The second key step in our preprocessing phase is the generation and selection of cutting planes. We have discussed various valid linear constraints and second order cone constraints that can be used to tighten the initial relaxation of LPCC in [49], and have shown the computational results of these valid constraints individually. As important as finding these cutting planes is the selection of the cuts that actually should be added to the initial LP relaxation. In this section, we will describe our detailed procedure to generate and select our cutting planes. Note that we will only add cutting planes at the root node, and perform the generation of each type of cut in rounds and in the following order:

- Disjunctive cuts and Simple cuts

- Bound cuts

- Linear cuts derived from second order cone constraints

We use the computational results with these cutting planes in [49] to guide the cut generation process. The details of generation and selection rule are described as follows.

input : the LP relaxation solution of the original LPCC:  $x^*$ ,  $y^*$ ,  $w^*$ , search depth parameter depth, search breadth parameter breadth output: recovered feasible LPCC solution or failed to recover the solution Initialization: Set binary vector  $z^* = 0$ ; for  $i \leftarrow 1$  to m do if  $y_i^* < w_i^*$  then  $z_i^* = 0;$ else  $z_i^* = 1;$  $\mathbf{end}$ Solve (4) to get  $FG-LPCC(z^*)$ ; if  $FG-LPCC(z^*) == 0$  then solve  $LPCC(z^*)$ , and **return** the optimal solution to  $LPCC(z^*)$ ;  $\mathbf{end}$ else let  $A(z^*)$  denote the set of binary vectors that are adjacent to  $z^*$ ; foreach  $z \in A(z^*)$  do solve (4) to get FG-LPCC(z); insert z into a sorted queue Q with nondecreasing order on FG-LPCC(z);  $\mathbf{end}$ Let  $r_b = 0;$ while Q is not empty and  $r_b \leq breadth \mathbf{do}$  $r_b = r_b + 1;$ pop the top element  $\bar{z}$  in Q, and delete this element from Q; let  $z = \overline{z}$  and  $r_d = 0$ ; while there exists any feasibility gap descent direction  $\Delta z$  for z and  $r_d \leq depth \operatorname{do}$ pick a feasibility gap descent direction  $\Delta z$ ;  $z = z + \Delta z;$  $r_d = r_d + 1;$  $\mathbf{end}$ if FG-LPCC(z) == 0 then solve LPCC(z), and **return** the optimal solution to LPCC(z); end  $\mathbf{end}$ end return feasibility recovery failed;

Procedure 1: Local search feasibility recovery process

```
input : the known valid upper bound of LPCC ub_{initial}, parameter searchGap_{min}
output: refined feasible LPCC solution or failed to refine the known feasible solution
Initialization: Set lb_{search} =optimal value of the LP relaxation of LPCC and ub_{search} = ub_{initial}; add lb_{search} \leq c^T x + d^T y \leq ub_{search} into (4);
while ub_{search} - lb_{search} > searchGap_{min} do
    solve LP relaxation of LPCC with constraints lb_{search} \leq c^T x + d^T y \leq ub_{search} ;
    apply Procedure 1 to recover a feasible solution;
    if recovery process succeed then
         update the refined feasible solution with recovered solution;
         update ub_{search} with the newly recovered solution;
        ub_{search} = (lb_{search} + ub_{search})/2;
    \mathbf{end}
    else
     lb_{search} = (lb_{search} + ub_{search})/2;
    end
\mathbf{end}
if refined feasible solution has been updated then
return refined feasible solution
\mathbf{end}
else
return feasibility refinement failed
end
```

**Procedure 2:** Refined local search feasibility recovery process

## 2.2.1 Disjunctive cuts and simple cuts

These cuts exploit the disjunctive constraints: for each *i*, either  $y_i \leq 0$  or  $w_i \leq 0$ . The solution to the LP relaxation typically violates a number of these disjunctions, and disjunctive cuts can either be generated by solving a supplemental linear program, or by examining the optimal tableau for the LP relaxation. Based on our computational experience, it seems that general disjunctive cuts and simple cuts [9,4,6] are the weakest cuts among our three type of cutting planes, but they are the cheapest to generate. Therefore we generate this type of cut first. The solving time of CPLEX for our test instances became worse when we added all of the generated disjunctive cuts or simple cuts to the root node even though the value of the initial LP relaxation was improved by these cuts, because the initial LP became too large. Moreover, due also to the high cost of generating general disjunctive cuts, we only generate  $\lfloor m/100 \rfloor$  rounds of general disjunctive cuts and for each round we only generate at most 3 general disjunctive cuts instead of generating disjunctive cuts for each violated complementarity constraint.

The values of  $y_i w_i$  in the optimal solution to the LP relaxation are sorted in nonascending order and we select complementarity constraints with index that corresponds to the largest three products. After each round of generating cuts, we will remove every cut whose corresponding slack variable is basic in the relaxed LP, in order is to keep the size of the relaxed LP small. After generating the general disjunctive cuts,  $\lfloor m/10 \rfloor$  rounds of simple cuts will be added. Since a simple cut is derived from the simplex tableau with almost no cost, we will generate simple cuts for every violated complementarity constraint in each round, and also remove every cut whose corresponding slack variable is basic in the relaxed LP after each round of generating cuts.

#### 2.2.2 Bound cuts

Upper bounds  $u_i^y$  and  $u_i^w$  on  $y_i$  and  $w_i$  can be used in the bound cut

$$u_i^w y_i + u_i^y w_i \le u_i^w u_i^y \tag{5}$$

for any pair of complementary variables  $y_i$  and  $w_i$ . Strengthening the upper bounds seems very important for the branch-and-bound routine of CPLEX for solving our instances, and the bound cuts also improve the initial lower bound dramatically. However, the major drawback of bound cuts is that they are very expensive to generate, especially when m, the number of complementarity constraints, is very large. Therefore, we will only compute bounds for at most 5 pairs of complementary variables, and the selection of these complementary variables is the same as the selection of complementarity constraints to generate disjunctive cuts. An upper bound  $u_i^y$  for  $y_i$  can be found by solving the linear program

$$u_{i}^{y} = \underset{(x,y,w)}{\operatorname{maximize}} y_{i}$$
(6)  
subject to  $Ax + By \ge b$   
 $x \ge 0$   
 $0 \le y \le u^{y}$   
 $0 \le q + Nx + My = w \le u^{w}$   
 $c^{T}x + d^{T}y \le ub$   
 $u_{j}^{w}y_{j} + u_{j}^{y}w_{j} \le u_{j}^{w}u_{j}^{y} \quad \forall j \text{ with known bounds } u_{j}^{y}, u_{j}^{w}$ 

where ub is a known upper bound on the optimal value of the LPCC. A similar LP can be constructed to get bounds on w.

We also investigated improving the bound cuts by splitting the variables. In particular, two versions of problem (6) could be solved, one with the additional constraint  $y_k = 0$  and the other with the additional constraint  $w_k = 0$ , for some index  $k \neq i$ . The maximum of the optimal values of these two problems could potentially improve on the initial upper bound. For our test instances, the additional computational work involved in computing these improved bounds did not improve the overall computational time, so this splitting is not included in our results.

## 2.2.3 Linear cuts from second order cone constraints

Based on the computational results of [49], cuts derived from a certain second order cone constraint can significantly improve the initial lower bound of our instances with relatively low generating cost compared to bound cuts when  $n \ll m$ , provided M is positive semidefinite. These cuts arise from linearizing the term  $y^T N x$ , using McCormick inequalities [48] to tighten the linearization, and handling the  $y^T M y$  term appropriately. Details can be found in [50]. The constraints can be tightened by refining bounds. We did not use these cuts in the computational results reported in this paper, because of difficulties with ensuring M was regarded as numerically positive semidefinite by CPLEX.

# 2.3 Overall Flow of the Preprocessor

The preprocessor consists of the following steps:

- 1. Apply the feasibility recovery routine to recover a feasible solution.
- 2. Generate  $\lfloor m/100 \rfloor$  rounds of general disjunctive cuts.
- 3. Generate  $\lfloor m/10 \rfloor$  rounds of simple cuts.
- 4. Apply 4 bound refinements and generate bound cuts.

We apply Procedures 1 and 2 as the default feasibility recovery procedure due to run time considerations. Other feasibility recovery procedures and refinements can also be invoked if required for solving special classes of problems. The number of rounds for generating each type of cutting plane can be modified by changing the parameter settings. The current setting is based on the computational experience in [49].

An additional preprocessing procedure undertaken at each node is the complementary variable fixing process, which is detailed in §3.3.

## **3** Branch-and-Bound Phase

The branch-and-bound routine needs to be invoked to solve the problem exactly if the initial LP relaxation is unbounded or the preprocessing phase is unable to close 100% of the gap for the bounded case. The branching is imposed on the complementarity constraint directly, and two subproblems (nodes) will be generated by enforcing either side of the pair of complementary variables to its lower bound zero. Just like a branch-and-bound based MIP solver, there are two key ingredients in our branch-and-bound routine: branching complementarity selection and node selection. Branching complementarity selection is the procedure to select the complementarity constraint to be branched on, and it is the same as the "variable selection" in mixed integer programming. In §3.1, we present our branching strategy which is based on the ideas of three classic branching rules and also some new proposed ideas designed for the LPCC problem. Node selection is the procedure to select the next subproblem from the node tree to be processed. In §3.2, we will present and compare different node selection strategies. Besides these two key ingredients, in §3.3 we will describe the node pre-solving procedure used in our algorithm to pre-process the nodes during the branch-and-bound process. The general branch-and-bound routine for handling the bounded case and unbounded case of LPCC are described in §4.1 and §4.2 respectively.

### 3.1 Branching Complementarity Selection

The branching rule is the key ingredient of any branch-and-bound algorithm. Good branching strategies are extremely important in practice for solving mixed integer programs, although currently there is no existing theoretical best branching strategy. We will first present three classic branching strategies for solving mixed integer programs that have been studied in the literature. The reader can refer to Linderoth and Savelsbergh [45], Fügenschuh and Martin [25] and Achterberg et al [2] for a comprehensive study of branch-and-bound strategies for mixed integer programming. We will present our branching strategy based on the ideas of these branching strategies. The computational results that compare various branching strategies will be shown in §5.

We first give some definitions related to our branching routine for the LPCC problem. For easy discussion, if the LP relaxation of the LPCC is unbounded below, we represent its lower bound as  $-\infty$ . Suppose that we have an LPCC problem Q and the set I is the index set of complementarity constraints. If the current solution to the LP relaxation of Q is not a feasible solution to LPCC (for the unbounded case, we consider an unbounded ray of the LP relaxation instead of solution to the LP relaxation), then

we can pick an index  $i \in I$  with  $y_i w_i > 0$  and obtain two subproblems (nodes): one by adding the constraint  $y_i \leq 0$  (named the left child node, denoted by  $Q_i^{i}$ ) and one by adding the constraint  $w_i \leq 0$ (named the right child node, denoted by  $Q_i^w$ ). We refer to this as branching on complementarity *i*. For the bounded case, if we denote the objective value of the LP relaxation of Q as  $c_Q$  and the objective value of the LP relaxation of its two child nodes as  $c_{Q_i^y}$  and  $c_{Q_i^w}$  respectively, then the objective value changes caused by branching on the *i*th complementarity are  $\Delta_i^y = c_{Q_i^y} - c_Q$  and  $\Delta_i^w = c_{Q_i^w} - c_Q$ . We usually use the improvement of objective value of the LP relaxation to measure the quality of branching on the *i*th complementarity. Our implementation supports fixing multiple complementarity constraints at one time, but by default we will only select to branch on one complementarity. Based on the results of testing our instances and the computational results of solving various MIP problems in the literature, multiple way branching is rarely better than two way branching.

The generic procedure for selecting the branching complementarity can be described in Procedure 3. The score function in Step 2 of this procedure needs to evaluate the two child nodes that could be

input : the LP relaxation solution of the current processing node Q or the unbounded ray to the LP relaxation if the LP relaxation is unbounded:  $x^*, y^*, w^*$ 

- **output:** the selected branching index  $i \in I$  of a complementarity constraint
- 1. Let  $\tilde{I} = \{j \in I \mid y_i^* w_i^* > 0\}$  denote the index set of violated
- complementarity constraints.
- 2. Compute a branching score  $s_j \in \mathbb{R}^+$  for all candidates  $j \in \tilde{I}$ . 3. Select the selected branching index  $i \in \tilde{I}$  with  $s_i = \max_{k \in \tilde{I}} \{s_k\}$ .

Return selected branching index i.

s

**Procedure 3:** Generic complementarity selection procedure

generated by the branching, and map these two effectiveness values onto a single score value. Different choices for the effectiveness values are given later. Suppose  $q^y$  and  $q^w$  are the effectiveness values of the two child nodes generated by a branching. In the literature, the score function usually has one of the following forms:

$$core(q^{y}, q^{w}) = (1 - \mu) \cdot \min\{q^{y}, q^{w}\} + \mu \cdot \max\{q^{y}, q^{w}\}$$
(7)

or

$$score(q^y, q^w) = \max\{q^y, \epsilon\} \cdot \max\{q^w, \epsilon\}$$
(8)

Here  $\mu$  is a number between 0 and 1, and it is usually an empirically determined constant or a dynamic parameter adjusted through the course of branching process. We chose  $\epsilon = 10^{-6}$  to enable the comparison when either  $q^y$  or  $q^w$  is zero. Based on the computational experience in [1], the product form is superior to the weighted sum form for solving MIP problems. Therefore, in our algorithm, we chose to use the product form to map the effectiveness values from two child nodes onto a single value.

In the following we will present three classic branching strategies for solving an MIP in terms of our branching on complementarity scheme: Strong Branching (apparently originally developed in the work leading up to [3]), Pseudocost Branching [12] and Inference Branching [1]. In fact, all of these branching routines are just variants of Procedure 3 with different score functions.

## 3.1.1 Strong branching

The idea of Strong Branching [3] is to test the branching candidates by temporarily enforcing either side of a complementarity constraint and solving the resulting LP relaxation to a certain level, then select the one that can lead to the largest lower bound improvement. Full Strong Branching will compute  $\Delta_i^y$  and  $\Delta_i^w$  for each branching complementarity candidate  $i \in I$ , and use the score  $(\Delta_i^y, \Delta_i^w)$  as the effectiveness values in the form of either (7) or (8) as its score function. Full Strong Branching can be seen as the locally best branching strategy in terms of lower bound improvement. However the computational cost of Full Strong Branching is very high, since in order to evaluate the score function for each complementarity candidate, we need to solve two resulting LP relaxations to optimality. There are usually two ways to speed up Full Strong Branching: one is to only test a subset of the candidate set instead of considering all the candidates, and another is to perform a limited number of simplex iterations and estimate the objective value change based on that. In our branch-and-bound algorithm, we have implemented the *Full* Strong Branching routine, and also we adopt the former idea to speed up the *Full Strong Branching*: as long as the objective value of LP relaxation of either side of the child nodes hits some threshold, we will select this branching candidate and exit the selection routine; we set the median value of the lower bound of unsolved nodes in the current search tree as this threshold.

A version of strong branching was used by Fischer and Pfetsch [22] in their branch-and-cut approach for problems with overlapping SOS1 constraints.

#### 3.1.2 Pseudocost branching

*Pseudocost Branching* [12] uses the branching history to estimate the two objective changes of the child nodes without actually solving them. In other words, *Pseudocost Branching* is a branching rule based on the historical performance of complementarity branching on complementarities which have already been branched. Let  $\varsigma_i^y$  and  $\varsigma_i^w$  be the objective gain per unit change at node Q after branching on complementarity i by enforcing  $y_i$  or  $w_i$  to zero, that is

$$\varsigma_i^y = \frac{\Delta_i^y}{y_i^*} \text{ and } \varsigma_i^w = \frac{\Delta_i^w}{w_i^*}$$

$$\tag{9}$$

where  $y_i^*$  and  $w_i^*$  are the violation of complementarity *i* corresponding to the LP relaxation solution of Q. Let  $\sigma_i^y$  denote the sum of  $\varsigma_i^y$  over all the processed nodes where complementarity *i* has been selected as the branching complementarity and resulting child node  $Q_i^y$  has been solved and was feasible. Let  $\eta_i^y$  denote the number of these problems, and define  $\sigma_i^w$  and  $\eta_i^w$  in the same way for the other side of the complementarity. Then the pseudocost of branching on complementarity *i* can be calculated as the arithmetic mean of objective gain per unit change:

$$\Psi_i^y = \frac{\sigma_i^y}{\eta_i^y} \text{ and } \Psi_i^w = \frac{\sigma_i^w}{\eta_i^w}$$
(10)

Therefore given the violated complementarity *i* corresponding to the LP relaxation of Q, it is reasonable to use  $\Psi_i^y \cdot y_i^*$  and  $\Psi_i^w \cdot w_i^*$  to estimate  $\Delta_i^y$  and  $\Delta_i^w$  respectively. We call the branching rule that uses the score function  $score(\Psi_i^y \cdot y_i^*, \Psi_i^w \cdot w_i^*)$  in step 2 of Procedure 3 as *Pseudocost Branching*. Notice that at the beginning of the branch-and-bound procedure, the pseudocost is uninitialized for all the complementarities. One way to handle a complementarity with an uninitialized pseudocost is to replace its pseudocost with the average of the pseudocosts of the complementarities are uninitialized. Applying strong branching to the nodes whose tree depth level is less than a given level is another way to initialized the pseudocosts. More recently, Achterberg et al [2] proposed a more general pseudocost initialization method, and named the corresponding branching rule as *Reliability Branching*. In our implementation, we include the pseudocost as part of our branching score, and we choose to apply strong branching to nodes whose tree depth level is less than pseudocost.

# 3.1.3 Inference branching

The branching decision of strong branching and pseudocost branching are both based on the change of objective value of the LP relaxation, while *Inference Branching* [1] is quite different from the above two branching strategies. *Inference Branching* checks the impact of branching on changing the bounds of other variables. As with pseudocosts, historical information is typically used to estimate the deductions on bounds of the variables, and the inference value can be calculated as the arithmetic mean of the number of bound deductions. The domain propagation process is a node pre-solving process to detect the bound change of the variables and is discussed in §3.3. In our implementation, we use a similar idea to inference branching: instead of evaluating the inference value, we estimate the *complementarity satisfaction level* after branching on a complementarity, leading to the quantity  $s_i^{SL}$  below.

# 3.1.4 Hybrid branching strategy for the LPCC (bounded case)

Our branching strategy for the bounded case combines the ideas of the above three classic branching strategies, and additionally we also include some new score values into our branching score function which are specialized for the LPCC problem.

In our implementation, the default branching strategy will apply the full strong branching strategy for the nodes whose depth level are no larger than 7. The reason for doing that is because it is usually quite important to make the right branching decision at the beginning, and also we can use strong branching to initialize the pseudocosts and another score value that we will propose next. For the nodes whose tree depth are larger than 7, we will use a weighted sum formula to combine four score values for each violated complementarity. Among these four score values, two of them are only based on the current node Q, and the other two are based on historical branching information. For the violated complementarity i, these four score values are listed as follows:

1.  $s_i^{VL}$ : score of Violation Level. Suppose  $y_i^*$  and  $w_i^*$  are the violation of complementarity *i* corresponding to the LP relaxation of Q, then we define

$$s_i^{VL} = \sqrt{y_i^* \cdot w_i^*}$$

2.  $s_i^{ED}$ : score of Euclidean Distances from the LP relaxation solution of Q to the two hyperplanes corresponding to  $y_i = 0$  and  $w_i = 0$ . Recall that since  $y_i^* \cdot w_i^* > 0$ , we can represent the complementary variables  $y_i$  and  $w_i$  with the non-basic variables in the optimal simplex tableau of Q

$$y_i = y_i^* - \sum_{j \in NB} a_j^{y_i} \xi_j \tag{11}$$

$$w_i = w_i^* - \sum_{j \in NB} a_j^{wi} \xi_j \tag{12}$$

We use the Euclidean distance from the LP relaxation solution to the two hyperplanes

$$\sum_{j \in NB} a_j^{yi} \xi_j = y_i^* \text{ and } \sum_{j \in NB} a_j^{wi} \xi_j = w_i^*$$

to define  $s_i^{ED}$  as follows:

$$s_i^{ED} = \sqrt{\frac{y_i^* \cdot w_i^*}{\sqrt{\|a^{yi}\| \cdot \|a^{wi}\|}}}$$

3.  $s_i^{PC}$ : score of Pseudo Cost. We use the following small modification to the pseudcost calculation of §3.1.2:

$$s_i^{PC} = \sqrt{\max\{\Psi_i^y \cdot y_i^*, \epsilon\} \cdot \max\{\Psi_i^w \cdot w_i^*, \epsilon\}}$$

4.  $s_i^{SL}$ : score of complementarity Satisfaction Level. We define the complementarity satisfaction level as the proportion of the satisfied complementarities corresponding to the LP relaxation solution of the child node after branching. Intuitively we want to select a branching complementarity that will lead to more satisfied complementarities. To estimate this complementarity satisfaction level, we collected the historical information to compute the average complementarity satisfaction level for both sides of the complementarity

$$\Phi_i^y = rac{\varphi_i^y}{\eta_i^y}$$
 and  $\Phi_i^w = rac{\varphi_i^w}{\eta_i^w}$ 

Here  $\varphi_i^y$  is the sum of the proportion of complementarity satisfaction levels over all the prior nodes, where complementarity *i* has been selected as the branching complementarity, and  $\eta_i^y$  is the total number of these nodes. We define  $\varphi_i^w$  and  $\eta_i^w$  to be the analogous value for the other side of complementarity. Then the score of the complementarity Satisfaction Level can be calculated as

$$s_i^{SL} = \sqrt{\Phi_i^y \cdot \Phi_i^w}$$

We scale the score vectors using their 2-norms, and the following formula is the branching score function that we used to evaluate the score for each violated complementarity:

$$s_{i} = \omega^{VL} \left( \frac{s_{i}^{VL}}{\|s^{VL}\|} \right) + \omega^{ED} \left( \frac{s_{i}^{ED}}{\|s^{ED}\|} \right) + \omega^{PC} \left( \frac{s_{i}^{PC}}{\|s^{PC}\|} \right) + \omega^{SL} \left( \frac{s_{i}^{SL}}{\|s^{SL}\|} \right)$$
(13)

By default, the weight is set as  $\omega^{VL} = 1$ ,  $\omega^{ED} = 0.5$ ,  $\omega^{PC} = 0.25$  and  $\omega^{SL} = 0.5$ . Note that setting different weights for each score value will lead to different branching behaviour. In §5.2, we will show the computational results of solving our LPCC instances with different weights of the score value.

## 3.1.5 Hybrid branching strategy for LPCC (unbounded case)

Our branching strategy for the unbounded case is slightly simpler than the one for the bounded case. We will still apply full strong branching to the nodes whose tree depth level is no larger than 7. However, for the remaining unbounded nodes we will only use  $s_i^{VL}$  as the branching score to make the branching decision.

#### 3.2 Node Selection

In addition to selecting which complementarity to branch on, another question is which subproblem (node) we should pick to process. There are two major criteria for selecting the next subproblem to be processed.

- 1. finding feasible LPCC solutions to improve the upper bound of the LPCC problem which leads to pruning the nodes by bounding, leading to a Depth First Search strategy.
- 2. improving the lower bound as fast as possible, leading to a Best-Bound strategy.

In our implementation of the branch-and-bound routine we use a **Best-Bound** strategy to select the next node to be processed, since we want to solve the problem to optimality as fast as possible. Notice that for the *Best-Bound*, it is possible that there are several nodes with the same lower bound. For that case, we will select the most recently generated node as the next node to be processed.

## 3.3 Node Pre-solving

The major task of our node pre-solving procedure is to tighten the domains of complementary variables  $y_i$  and  $w_i$  and try to fix the complementary variables. In order to facilitate the discussion, here we can assume that each  $\xi_i$  in (11) and (12) is a non-negative variable with zero lower bound. Therefore we have the following result: if  $a_j^{y_i} \leq 0, \forall j \in NB$ , then we have  $y_i \geq \hat{y}_i$ , and therefore  $w_i = 0$ ; if  $a_j^{w_i} \leq 0, \forall j \in NB$ , then we have  $w_i \geq \hat{w}_i$ , and therefore  $y_i = 0$ . This complementary variable fixing check is performed before we branch on the complementarity constraint.

## 4 General Scheme of the Branch-and-Cut Algorithm for Solving LPCC

The preprocessing routines are only invoked if the the initial LP relaxation of the LPCC has a bounded optimal value; we refer to this as the "bounded case". If the initial relaxation does not have a finite optimal value then we are in the "unbounded case". For the bounded case, the preprocessing procedure is applied first to tighten the initial LP relaxation, then the branch-and-bound routine is invoked to solve the LPCC to optimality; for the unbounded case, we will only apply the branch-and-bound routine, which gives unbounded nodes higher priority than bounded nodes. A flow diagram of the overall algorithm is given in Figure 1. The initialization step 0 sets the upper bound  $\bar{z} = +\infty$ , the lower bound  $\underline{z} = -\infty$ , the unbounded node list  $\bar{L} = \emptyset$ , and the bounded node list  $L = \emptyset$ . If the LP relaxation of the initial problem is feasible then the initial problem is added to L or  $\bar{L}$  in box 1, as appropriate. Boxes 2, 4, 6, 8, 10, 12, and 14 corresponding to the bounded case are the subject of §4.1, with the unbounded case boxes 3, 5, 7, 9, and 11 explained in §4.2. Box 13 is discussed in §4.3.

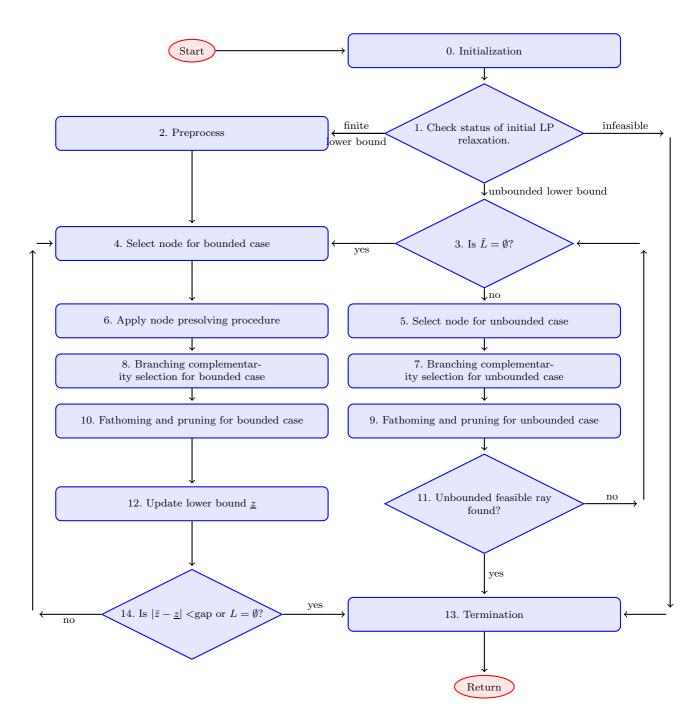


Fig. 1 Flow chart of branch-and-bound procedure

# 4.1 Overall Flow of Branch-and-Bound for LPCC (Bounded Case)

In the bounded case, the algorithm is quite similar to the branch-and-bound routine for a mixed integer program. If it is determined in box 1 that the initial LP relaxation is bounded then we implement a more detailed preprocessing step in box 2, as discussed in §2. In box 4, we apply *Best-Bound* to pick the next node  $LPCC^i$  from L to be processed and delete  $LPCC^i$  from L. The node presolving procedure from §3.3 is implemented in box 6. The branching strategy of Section 3.1.4 is used in box 8 to select branching complementarity j. Fathoming and pruning is performed in box 10 as follows:

Fathoming and pruning: Generate two child nodes by enforcing either  $y_j = 0$  or  $w_j = 0$  and solve LP relaxations. For each child node:

- 1. If LP relaxation solution is feasible in LPCC with objective  $z^*$  then delete child node. Set  $\bar{z} \leftarrow \min\{\bar{z}, z^*\}$ .
- 2. If LP relaxation is feasible with objective  $z^* < \overline{z}$  then set the lower bound of child node as  $z^*$  and add child node to L.
- 3. If LP relaxation feasible with objective  $z^* \geq \overline{z}$  or infeasible then delete child node.

The lower bound is updated in box 12. The procedure is terminated in box 14 if there are no more nodes in the set L or if the gap between the upper and lower bound is sufficiently small.

4.2 Overall Flow of Branch-and-Bound for LPCC (Unbounded Case)

The branch-and-bound routines for solving mixed integer programs in existing MIP solvers like CPLEX usually assume the initial LP relaxation is bounded below. Even if the initial LP relaxation is unbounded, it is still treated as bounded below by adding an objective lower bound constraint with a very large negative number  $(-10^{20})$  as its lower bound. However, our branch-and-bound routine for handling the unbounded case of the LPCC is quite different. If the LP relaxation of a node is unbounded, we will treat this node as an unbounded node and add it to the unbounded node list. If the unbounded node list is non-empty, our branch-and-bound routine will always process a node in the unbounded node list first. Notice that when we find an unbounded ray that satisfies all the complementarities, we need to check whether this is a feasible ray to the LPCC. The LPCC is feasible with unbounded objective value if and only if we find an unbounded feasible ray to the LPCC.

If the set  $\overline{L}$  of unbounded nodes is empty in box 3 then we return to the bounded case in box 4, constructing an appropriate lower bound  $\underline{z}$ . In box 5, we select the node  $LPCC^i$  that is the most recently generated from  $\overline{L}$  to be processed and delete  $LPCC^i$  from  $\overline{L}$ . The branching strategy of Section 3.1.5 is used in box 7 to select branching complementarity j. Fathoming and pruning for an unbounded node is performed in box 9 as follows:

Fathoming and pruning: Generate two child nodes by enforcing either  $y_j = 0$  or  $w_j = 0$  and solve LP relaxations. For each child node:

- 1. If LP relaxation solution is feasible in LPCC with objective  $z^*$  then delete child node. Set  $\bar{z} \leftarrow \min\{\bar{z}, z^*\}$ .
- 2. If LP relaxation is feasible with objective  $z^* < \overline{z}$  then set the lower bound of child node as  $z^*$  and add child node to the bounded node list L.
- 3. If LP relaxation feasible with objective  $z^* \geq \overline{z}$  or infeasible then delete child node.
- 4. If LP relaxation is unbounded and the unbounded ray is not a feasible ray to LPCC then add this child node to the unbounded node list  $\bar{L}$ .
- 5. If LP relaxation is unbounded and the piece of LPCC corresponding to that ray is feasible then the LPCC is unbounded.

If an unbounded piece is found in box 11 then the algorithm can be terminated; otherwise we loop back to box 3.

## 4.3 The Complete Overall Scheme

A flow chart of the algorithm is exhibited in Figure 1. Each of the three possible problem states can be returned in the termination box 13. If an unbounded feasible ray to the LPCC is found then the LPCC is feasible with unbounded objective value. If the LPCC is not unbounded and an LPCC feasible solution is found then the LPCC attains a finite optimal solution with optimal objective  $\bar{z}$ . Otherwise, the problem is infeasible.

## **5** Computational Results

In this section, we will present the computational results of using our proposed branch-and-cut algorithm to solve various LPCC instances. All procedures and algorithms are developed in the C language with the CPLEX callable library, and all LPs and convex quadratic constraint programs are solved using

1	m	rankM	Average gap	Optimal found out of 10
	100	30	0.09%	5
	100	60	0.22%	2
	150	30	0.0~%	10
	150	100	0.06%	3
	200	30	0.0~%	10
	200	120	0.07%	2

**Table 1** Average Computational Results of Feasibility Recovery with n = 2, k = 20. The column "Average gap" iscalculated as  $\frac{LB_{recovered} - LPCC_{opt}}{LPCC_{opt}}$ . Detailed results can be found in Table 5 in Appendix A.

CPLEX 12.6.2. We implement our algorithm through the addition of callback routines to CPLEX. As an alternative to our approach, CPLEX allows the modeling of complementarity constraints through the use of *indicator constraints*; we compare the computational performance of our algorithm with that of using default CPLEX 12.6.2 to solve indicator constraint formulations of these LPCC instances, with our preprocessor used for both approaches. Except for a few preliminary tests discussed in §5.2, all the computational testing is performed on a Mac Pro with 6 dual processor Intel Xeon E5 cores and 16GB of memory. Our branch-and-cut routine uses just one thread, while the default CPLEX 12.6.2 indicator constraint formulation can use all 12 available threads. The relative gap for optimality is  $10^{-6}$ , here the relative gap is defined as  $\frac{upperbound - lowerbound}{max(1, |lowerbound|)}$ . This is smaller than CPLEX's default MIP optimality tolerance and larger than its default LP tolerance. The tolerance of complementarity is  $10^{-6}$ , i.e., either  $y_i$  or  $w_i$  for i = 1, ..., m should be less than  $10^{-6}$  for any feasible LPCC solution. All runtimes are reported in seconds.

We used three sets of test instances. The first set consists of 60 LPCC instances with n = 2 and between 100 and 200 complementarities. The generation scheme for these problems and computational results can be found in Appendix A, with the results discussed in sections 5.1 and 5.2. The second set of test instances are LPCC formulations of bilevel programs, where the lower level problem is a convex quadratic program; the formulation and results are presented in Section 5.3, with more extensive results in Appendix B. The final set of results in Section 5.4 are for inverse quadratic programming problems, with detailed results in Appendix C.

Source code and test instances can be found online at https://github.com/mitchjrpi/LPCCbnc Also included with the source code is a Makefile. A user needs to have access to CPLEX in order to be able to compile the code. Generators for the bilevel and inverse QP problems can be found on the website; the generator uses AMPL to construct the instances.

## 5.1 Computational Results of the Feasibility Recovery Process

We will first apply the local search feasibility recovery process (procedure 1); if this procedure successfully recovers a feasible solution, then the refinement procedure (procedure 2) will be applied to refine that feasible solution. We set the depth parameter as 5 and breadth parameter as m, i.e. the number of complementarities, in procedure 1. Table 1 summarizes the feasibility recovery result of the 60 LPCC instances. The computational results show that our proposed feasibility recovery procedures can successfully recover a feasible solution for all of the 60 LPCC instances with very good quality. For most instances, the recovered feasible solution is in fact an optimal solution. Note that as m increases, the feasibility recovery processing time increases as well. Therefore in practice, as a preprocessing procedure, we need to control the depth and breadth parameters in procedure 2 to reduce the time spent on the feasibility recovery procedure.

#### 5.2 Computational Results of Branch-and-Cut Algorithm

In this section, we will show the computational results of using our proposed branch-and-cut algorithm to solve the 60 LPCC instances with finite global optimal values from Appendix A.

m	$Time_{R_1}$	$Time_{R_2}$	$Time_{R_3}$	$Time_{R_4}$	$Time_{CPLEX}$
	(sec)	(sec)	(sec)	(sec)	(sec)
100	19.185	18.790	18.805	18.917	38.021
150	75.213	73.623	76.071	74.285	1688.160
200	308.293	296.663	287.232	291.057	5043.017

Table 2 Comparison of geometric means of solving time, using our four different branching rules and using default CPLEX

ſ	m	$Node_{R_1}$	$Node_{R_2}$	$Node_{R_3}$	$Node_{R_4}$	$Node_{CPLEX}$
	100	213	215	212	196	39114
	150	831	863	848	757	1078311
	200	3408	3301	3161	2837	1494577

 Table 3 Comparison of geometric means of number of nodes in branch-and-cut tree, using our four different branching rules and using default CPLEX

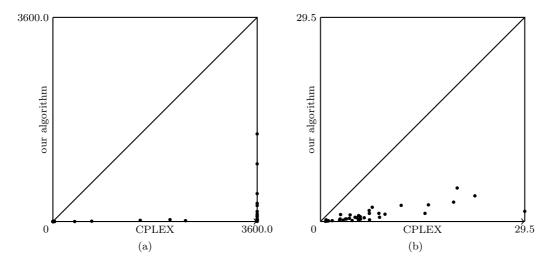


Fig. 2 Scatter plots for CPU time (seconds) for solution of LPCCs. Horizontal axis is time for the default CPLEX indicator constraint solver, vertical axis is time for our branch and cut algorithm. Processing times excluded. (a) All 60 instances. (b) 37 LPCCs where default CPLEX MIP required no more than 30 seconds.

We conducted preliminary experiments with 4 different weight settings of increasing sophistication to choose a score function (13):

 $\begin{array}{l} R_1: \ \omega^{VL} = 1, \ \omega^{ED} = 0, \ \omega^{PC} = 0 \ \text{and} \ \omega^{SL} = 0; \\ R_2: \ \omega^{VL} = 1, \ \omega^{ED} = 0.5, \ \omega^{PC} = 0 \ \text{and} \ \omega^{SL} = 0.5; \\ R_3: \ \omega^{VL} = 1, \ \omega^{ED} = 0.5, \ \omega^{PC} = 0.25 \ \text{and} \ \omega^{SL} = 0.5; \\ R_4: \ \omega^{VL} = 1, \ \omega^{ED} = 0.5, \ \omega^{PC} = 0.25 \ \text{and} \ \omega^{SL} = 0.5 \ \text{and} \ \text{apply strong branching rule to the node whose} \\ \text{tree depth is less or equal to 7.} \end{array}$ 

These results were obtained using CPLEX 11.4 using a single core of AMD Phenom II X4 955 CPU @ 3.2GHZ, 4GB memory and are contained in Tables 2 and 3. All four rules required far fewer nodes than default CPLEX. Based on these results,  $R_4$  is the best branching rule in terms of the number of nodes. Since in terms of solving time, these 4 routines are quite close, we chose  $R_4$  as our default branch-and-bound routine.

All remaining results in the paper were obtained using CPLEX 12.6.2 with detailed results contained in Table 6 in Appendix A. A scatter plot of the CPU time for solving the instances is given in Figure 2. Performance profiles [15] are given in Figure 3. The preprocessing times have been excluded from these plots. All the LPCC instances can be solved by our algorithm within thirty minutes, with 90% of them (54/60) solved within 150 seconds. Each instance requires considerably less processing time with our algorithm than with default CPLEX. Notice that default CPLEX is only able to solve 42 of the 60 instances within 3600 seconds. In particular, it is unable to solve 11 of our 20 LPCC instances when m = 200 within this time limit.

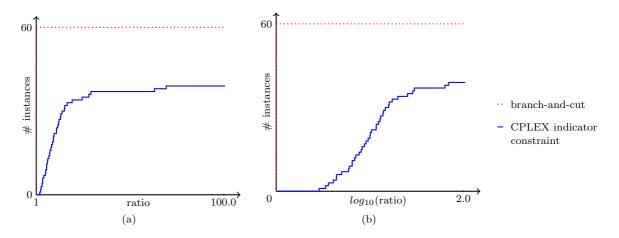


Fig. 3 Performance profile for CPU time (seconds) for solution of 60 LPCCs (preprocessing time excluded). Vertical axis is the number of instances. Horizontal axis is ratio of time required by the given algorithm to the time required by the better algorithm. (a) Linear scale. (b) Log scale.

The determination of a valid disjunctive cut or bound cut requires the solution of a linear programming problem. The parameter choices given in §2.2 result in 0.3m disjunctive cuts, approximately 5m simple cuts, and 15 bound cuts for each instance. We also experimented with not adding cutting planes in the preprocessor, in which case both codes performed slightly worse for the larger instances (a difference of perhaps 10% in average runtime for our branch-and-cut code).

# 5.3 Bilevel Test Problems

We further tested our algorithm on bilevel problems of the form

$$\min_{x,v} \quad c^T x + d^T v 
\text{subject to } Ax + Bv \ge b 
\quad 0 \le \quad v \le u 
\quad x \in \operatorname{argmin}_x \{ \frac{1}{2} x^T Q x + v^T x : Hx \ge g, x \ge 0 \}$$
(14)

where Q is positive semidefinite. The variables v are first stage variables, with the second stage variables x chosen to optimize a convex quadratic subproblem that depends on v. Both sets of variables appear in the linear objective. In addition, the first and second stage variables must satisfy the linking constraint  $Ax + Bv \ge b$ . By introducing KKT multipliers y and  $\lambda$  for the constraints in the subproblem, we can model this problem equivalently as the LPCC

$$\min_{x,v,y,\lambda,w} c^T x + d^T v \\ \text{subject to} \quad \begin{array}{l} Ax + Bv & \geq b \\ Qx + v - H^T y - \lambda &= 0 \\ 0 \leq v & \leq u \\ 0 \leq \lambda \perp x & \geq 0 \\ 0 \leq y \perp w := Hx - g \geq 0, \end{array}$$

a problem equivalent to one in our standard form (1). The relationship between the dimensions in (1) and the dimensions of the variables and constraints in (14) is as follows:

	Dimensions
(1)	(14)
m	$\operatorname{dimension}(g) + \operatorname{dimension}(v)$
n	$2 \times \operatorname{dimension}(v)$
k	dimension(b) + $3 \times \operatorname{dimension}(v)$

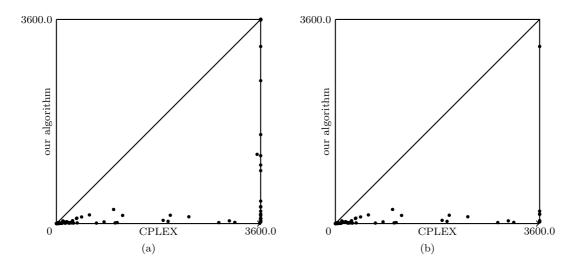


Fig. 4 Scatter plots for CPU time (seconds) for solution of LPCCs based on bilevel instances Horizontal axis is time for default CPLEX indicator constraint solver, vertical axis is time for our branch-and-cut solver. Processing times excluded. (a) All 90 instances. (b) 63 instances with n = 50.

1		Dimensi	on of $g$		Dimensi	on of b	Rank of $Q$			
	dim	time	# instances	dim	time	# instances	rank	time	# instances	
	50	12.76	16	25	39.98	16	25	19.94	31	
	100	30.29	15	50	51.06	16	50	161.66	32	
	150	41.91	16	75	15.11	16				
	200	266.45	16	100	259.57	15				

**Table 4** Average performance on bilevel instances with 50 first stage variables. Each column contains results from all 63 instances. Each average is taken over instances where the other parameters are varied.

Thus, the number of complementarity constraints is equal to the sum of the dimensions of g and v.

In our experiments, all parameters in b, c, d, g, A, B, and H were uniformly generated in the interval (0, 1). The matrix Q was equal to the matrix product  $LL^T$ , where the number of columns in L is equal to the required rank of Q and each entry in L is chosen uniformly from the interval (-1, 1). Each entry of u was equal to 1. Repeated problem dimensions in the table correspond to different instances. The dimension of g varied from 50 to 200, the dimension of v and x varied from 50 to 100, the number of complementarity constraints varied from 100 to 250, the dimension of b varied from 25 to 100, and the rank of Q varied between 0.5 of the dimension of v and the dimension of v. Problem data for the 90 bilevel test instances can be found in Tables 7 and 8.

We gave each algorithm a time limit of 3600 seconds in addition to the preprocessing time. Detailed performance data can be found in Tables 9 and 10. Our algorithm was able to solve all 63 instances with dimension of v equal to 50, 16/24 of the instances with the dimension of v equal to 75, and 3/3 of the instances with the dimension of v equal to 100. The corresponding numbers for the default CPLEX indicator constraint code were 56/63, 2/24, and 1/3. Our algorithm was considerably faster than default CPLEX indicator constraint code on every instance. Further, it had a smaller final gap than default CPLEX indicator constraint code for each instance where neither code could solve the problem. There was no instance that could be solved by default CPLEX indicator constraint code which could not also be solved by our algorithm. A scatter plot of the CPU time for solving the instances (ignoring the common preprocessing time) is given in Figure 4 and a performance profile is in Figure 5.

The instances become more difficult as the dimensions of v, b, and g increase, as might be expected. The instances also become more difficult as the rank of Q increases. Table 4 contains averages of solution times over these different parameters for the instances with the dimension of v equal to 50.

The parameter choices given in  $\S2.2$  result in 0.3m disjunctive cuts, approximately 3m simple cuts, and 15 bound cuts for each instance. Also as in  $\S5.2$ , we experimented with not adding cutting planes in the preprocessor. Both codes performed similarly to their respective performance with the preprocessor.

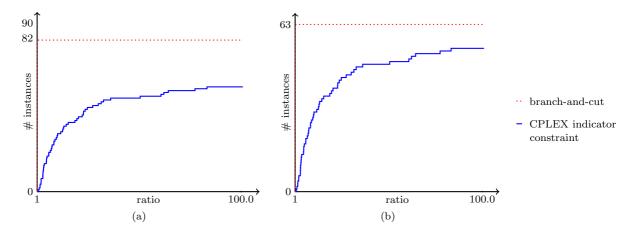


Fig. 5 Performance profile with linear scale for CPU time (seconds) for solution of LPCCs based on bilevel instances (preprocessing time excluded). Vertical axis is the number of instances. Horizontal axis is ratio of time required by the given algorithm to the time required by the better algorithm. (a) All 90 instances. (b) 63 instances with n = 50.

Thus, based on the results in this section and §5.2, our default implementation is to generate cutting planes in the preprocessor.

## 5.4 Inverse Quadratic Programs

Jara-Moroni et al. [37] presented a DC method for finding local optima for LPCCs arising from inverse quadratic programs [33]. The problem of interest has the form

$$\min_{x,b,c} ||(x,b,c) - (\bar{x},\bar{b},\bar{c})||_{1}$$
s.t.  $x \in \operatorname{argmin}_{y} \{ \frac{1}{2} y^{T} Q y + c^{T} y : A y \ge b \}$ 

$$(x,b,c) \in P$$

$$(15)$$

where  $\bar{x}$ ,  $\bar{b}$ , and  $\bar{c}$  are observations of the parameters and solution of a quadratic program and P is a polyhedron. The objective is to find (x, b, c) close to the observed values where x does solve the lower level quadratic program. In our computational testing, we varied the number of rows  $\tilde{m}$  and columns  $\tilde{n}$  of A between 100 and 400 and between 5 and 90, respectively; the dimensions of all other vectors and matrices are determined by the dimensions of A. When the matrix Q is positive definite, the inverse QP is equivalent to the following LPCC:

$$\min_{x,b,c,z^{x},z^{b},z^{c},\lambda} \mathbf{1}^{T} z^{x} + \mathbf{1}^{T} z^{b} + \mathbf{1}^{T} z^{c} 
s.t. \qquad Qx + c - A^{T} \lambda = 0 
\qquad x + z^{x} \ge \bar{x}, -x + z^{x} \ge -\bar{x} 
\qquad b + z^{b} \ge \bar{b}, -b + z^{b} \ge -\bar{b} 
\qquad c + z^{c} \ge \bar{c}, -c + z^{c} \ge -\bar{c} 
\qquad (x,b,c) \in P 
\qquad 0 \le \lambda \perp w := Ax - b \ge 0$$
(16)

where  $\lambda$  is the vector of KKT variables for the inner QP, the variables  $z^x$ ,  $z_b$ ,  $z^c$  are used to represent the  $L_1$  objective function in (15), and **1** represents a vector of ones of an appropriate dimension.

The instances in [37] were generated in MATLAB, whereas our instances were generated using AMPL. Nonetheless, we closely followed their procedures except for the generation of Q. Our matrix  $Q \in \mathbb{R}^{\tilde{n} \times \tilde{n}}$  was formed as the product  $MM^T$ , where  $M \in \mathbb{R}^{\tilde{n} \times \tilde{n}}$  was a square matrix with exactly three nonzeroes per row, with diagonal entries uniformly distributed between 0.5 and 1 and two off-diagonal entries uniformly distributed between 0 and 1; this results in a positive definite matrix Q, with about 9 entries per row on average (similar to the number of nonzeroes in a row of Q from [37]). Other parameters were generated as in [37]: the matrix  $A \in \mathbb{R}^{\tilde{m} \times \tilde{n}}$  has an average of approximately 10 nonzero entries per row which are uniformly distributed between 0 and 1; a vector  $\tilde{x} \in \mathbb{R}^{\tilde{n}}$  has components distributed as Normal(0,1);

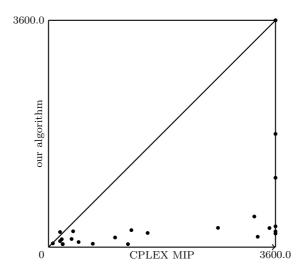


Fig. 6 Scatter plots for CPU time (seconds) for solution of inverse quadratic programs. Horizontal axis is time for CPLEX MIP called from AMPL and run on a single thread, vertical axis is time for our branch and cut algorithm.

vectors  $\hat{\lambda} \in \mathbb{R}^{\tilde{m}}$  and  $\hat{w} \in \mathbb{R}^{\tilde{m}}$  have components uniformly distributed between 0 and 10; a binary vector  $v \in \mathbb{B}^{\tilde{m}}$  is generated and  $\tilde{\lambda} \in \mathbb{R}^{\tilde{m}}$  and  $\tilde{w} \in \mathbb{R}^{\tilde{m}}$  are constructed as the Hadamard products  $\tilde{\lambda} := \hat{\lambda} \bullet v$  and  $\tilde{w} := \hat{w} \bullet (\mathbf{1} - v)$ ; vectors  $\tilde{b} \in \mathbb{R}^{\tilde{m}}$  and  $\tilde{c} \in \mathbb{R}^{\tilde{n}}$  are defined as  $\tilde{b} := A\tilde{x} - \tilde{w}$  and  $\tilde{c} = A^T\tilde{\lambda} - Q\tilde{x}$ ; vectors  $\bar{x} \in \mathbb{R}^{\tilde{n}}$ ,  $\bar{b} \in \mathbb{R}^{\tilde{m}}$ , and  $\bar{c} \in \mathbb{R}^{\tilde{n}}$  are obtained by perturbing  $\tilde{x}$ ,  $\tilde{b}$ , and  $\tilde{c}$  respectively, using Normal (0,1) noise; the polyhedron P is constructed as a box using simple bounds  $-u^x \mathbf{1} \le x \le u^x \mathbf{1}, -u^b \mathbf{1} \le b \le u^b \mathbf{1}, -u^c \mathbf{1} \le c \le u^c \mathbf{1}$  with  $u^x = 10 \max\{|\tilde{x}_i|\}, u^b = 10 \max\{|\tilde{b}_i|\}, u^c = 10 \max\{|\tilde{c}_i|\}$ ; finally, upper bounds are also imposed on  $\lambda$  with  $u^{\lambda} = 10 \max\{|\tilde{\lambda}_i|\}$ . The point  $(\tilde{x}, \tilde{b}, \tilde{c})$  with  $\tilde{\lambda}$  is feasible in the resulting problem instances of (16).

It is easy to generate explicit upper bounds on w = Ax - b from the upper bounds on x and b. Also, explicit upper bounds on  $\lambda$  are imposed following [37]. Thus, this problem can be formulated directly as a mixed integer program of the form (2). Because of this observation, our comparisons in this section are somewhat different from the previous experiments. In particular, we make the following two changes:

- Since bounds are already available, we do not use the cutting plane generation features of the preprocessor.
- We compare our LPCC branch-and-cut code with the CPLEX MIP solver invoked from AMPL, run with a single thread.

Our testbed consisted of 5 sets of 5 instances:  $(\tilde{m}, \tilde{n})$  equal to (100,75), (120,90), (150,20), (200,15), and (400,5). A scatter plot of the results can be found in Figure 6 and performance profiles can be found in Figure 7. Detailed computational results are contained in the Appendix, in Table 11. Our algorithm was able to solve 23 of the 25 instances within the 3600 second time limit; the corresponding figure for CPLEX was 18 out of 25. There was only one instance where CPLEX outperformed our code. Our algorithm solved 20 of the 25 instances within 360 seconds, while CPLEX only solved 6 of the instances within this time window.

## 6 Conclusions

The optimal solution to a linear program with complementarity constraints can in principle be found directly using CPLEX. However, far better performance can often be obtained by adding good cutting planes, by incorporating a specialized feasibility recovery routine, and especially by designing good branching routines. Our computational results demonstrate that our code is at least an order of magnitude faster than a default version of CPLEX, at least for our test set of instances. It is able to solve instances with up to 400 complementarity constraints in reasonable amounts of time, and can reliably solve instances with 100 complementarity constraints in less than a minute.

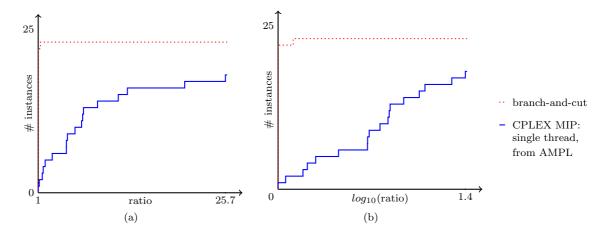


Fig. 7 Performance profile for CPU time (seconds) for solution of 25 inverse quadratic programs. Vertical axis is the number of instances. Horizontal axis is ratio of time required by the given algorithm to the time required by the better algorithm. Time limit 3600 seconds. (a) Linear scale. (b) Log scale.

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input : n,m,k,rankM,denseoutput: vector c,d,b,q; matrix A,B,N,M1: generate n dimension vector  $\bar{x}$  with value between 0 and 10, integer; 2: generate m dimension vector  $\bar{y}$  with value between 0 and 10, integer if index  $<\frac{m}{3}$ ; 0 otherwise; 3: generate n dimension vector c with value between 0 and 10, integer; 4: generate m dimension vector d with value between 0 and 10, integer; 5: generate  $k \times n$  matrix A with value between -5 and 6, integer, and the matrix density is dense; 6: generate  $k \times m$  matrix N with value between -5 and 6, integer, and the matrix density is dense; 7: generate  $m \times n$  matrix N with value between -5 and 6, integer, and the matrix density is dense; 8: generate  $m \times n$  matrix N with value between -5 and 6, integer, and the matrix density is dense; 8: generate  $m \times n$  matrix  $\Delta M$  with value between -2 and 2, integer; Let  $m \times m$  matrix  $M = LL^T + \Delta M - \Delta M^T$ ; 9: generate k dimension vector  $\Delta b$  with value between 1 and 11, integer; let k dimension vector  $b = A\bar{x} + B\bar{y} - \Delta b$ ; 10: generate m dimension vector  $\Delta q$  with value 0 if index  $<\frac{2m}{3}$ ; integer between 1 and 11 otherwise; let m

dimension vector  $q = -N\bar{x} - M\bar{y} + \Delta q;$ 

Procedure 4: LPCC instances generator

# A LPCC Test Instances

In order to test the effectiveness of different type of valid constraints, a series of LPCC instances was randomly generated, and Procedure 4 gives a detailed description of the generator.

Remark 1 In the initialization step of the procedure, n is the dimension of x variable; m is the dimension of y variable; k is the dimension of b; rankM is the rank of matrix M; dense is the density of generated matrices; we assume all instances have the non-negativity constraint  $x \ge 0$  which are not included in the constraint  $Ax + By \ge b$ ; step 1 and step 2 are used to generate a feasible LPCC solution; step 8 is to generate matrix M to be a non-symmetric positive semidefinite matrix with rank rankM.

We generated 60 LPCC instances with 100, 150, 200 complementaries, 20 instances of each size, and with the same parameter, we randomly generated 5 instances. For CPLEX solving LPCC instances, we used indicator constraints in CPLEX C callable library [35] to formulate the complementarity constraints, and the CPLEX setting is default. The time limit for CPLEX is 3600 seconds. Notice that default CPLEX is unable to solve most of our LPCC instances when m = 200 within 3600 seconds. Table 5 contains objective function value information for the 60 instances, including the effectiveness of the preprocessing routines.

Table 6 contains performance data.

# **B** Bilevel Test Instances

Our code solved all 63 of the instances with dimension of v equal to 50 and 18/35 of the larger instances. With extended time, default CPLEX was able to solve all but one problem with n = 50; it still has a gap of 16.56% for problem 60 after more than 7200 seconds of wall clock time and 47304 seconds of processor time. It solved just 6/35 of the larger instances. Run time information can be found in Tables 9 and 10.

# C Inverse QP Instances

Computational results on 25 inverse QP instances can be found in Table 11. For each set of 5 instances, the average CPU time is listed if all the instances were solved or the number of solved instances is noted.

					Optimal	LP	Preprocessee	d bounds	Relative	e gaps (per	centages)
	#	m	$\operatorname{rank}$	dense	Value	relaxation	lower			UB-opt	% closed
$  \begin{array}{ccccccccccccccccccccccccccccccccccc$	1	100	30	70	769.911528	629.002874	669.22439	770.287	13.13	0.05	
$  \begin{array}{ccccccccccccccccccccccccccccccccccc$						650.929154				0.35	
						627.332027	657.571025				
		100		70	543	531.188245	539.856029	544.497	0.85	0.28	
$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	5	100	30	70	930		896.57115	930.917		0.10	
							588.868287				
						425.717966	459.942655				
	8	100	30	20	771	687.744893	745.909078	771	3.25	0.00	
	-						620.866694			0.00	
	10	100				705.051229	729.547378			0.00	
	11			70		606.45432	609.833638		2.03	1.66	
	12	100	60	70	686.130259		675.208522				
	13	100	60	70		722.033536		734			
	14	100	60	70	665.868588	657.703283	661.460391		0.68		
$  \begin{array}{ccccccccccccccccccccccccccccccccccc$	15	100	60	70	984.588193		855.932906	986			
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	16	100	60	20	691	629.620621	664.054558	691	3.90	0.00	43.90
$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	17	100	60	20	666.995818	631.110603	655.171515	667	1.77	0.00	32.95
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	18	100	60	20	756.780603	725.103749	746.527684	758	1.52	0.16	32.37
$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	19	100	60	20	763	626.529227	722.010148	763.971	5.50	0.13	30.04
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$					532.218697	521.894551					
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	21	150	30	70	1029	946.929565	1010.002422	1029	1.85	0.00	23.15
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	22	150	30	70	1160	1075.719667	1143.215912	1160	1.45	0.00	19.91
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	23	150	30	70	965	929.722695	957.060812	965	0.82	0.00	22.51
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	24	150	30	70	1242	1170.744571	1232.634488	1242	0.75	0.00	13.14
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	25	150	30	70		1013.045865	1063.947928	1149	7.40	0.00	62.56
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	26	150	30	20	822.333333	790.161133	813.932095	822.333	1.02	0.00	26.11
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	27	150	30	20	1046	991.351886	1039.478766	1046	0.62	0.00	11.93
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	28	150	30	20	922	851.085225	899.489258	922	2.44	0.00	31.74
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	29	150	30	20	992	855.028214	921.051941	992	7.15	0.00	51.80
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	30	150	30	20	848	729.617101	775.254605	848	8.58	0.00	61.45
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	31	150	100	70	1377.072388	1263.798462	1344.135656	1377.072	2.39	0.00	29.08
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	32	150	100	70	837	833.238632	835.993215	837	0.12	0.00	26.77
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	33	150	100	70		912.297933	951.089989	972.804	2.23	0.00	35.86
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	34	150	100	70		1206.833191	1238.300018	1261.188	1.82	0.05	41.45
$ \begin{array}{cccccccccccccccccccccccccccccccccccc$		150	100	70	1087.08492	1040.170448	1077.111477	1089		0.18	
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	36	150	100	20	921.273479	893.518557	904.290053	923		0.19	61.19
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	37	150	100	20	923.772654	774.71571		925			29.59
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	38	150	100	20							
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	39	150	100	20	879.582356			879.605		0.00	41.07
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	40	150		20	1158.383138	1063.017814	1119.548217	1158.432	3.36	0.00	40.72
$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	41	200	30	70	1580	1098.044624	1196.5995	1580	24.27	0.00	79.55
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$											
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57200120201224.7646831100.945211188.73487412263.040.1029.1058200120201145.9697921132.9963191140.09391711470.600.0945.29592001202014261399.2252511415.851591429.3640.950.2437.9060200120201371.9019591340.7844151358.03524413721.020.0144.56											
58         200         120         20         1145.969792         1132.996319         1140.093917         1147         0.60         0.09         45.29           59         200         120         20         1426         1399.225251         1415.85159         1429.364         0.95         0.24         37.90           60         200         120         20         1371.901959         1340.784415         1358.035244         1372         1.02         0.01         44.56											
592001202014261399.2252511415.851591429.3640.950.2437.9060200120201371.9019591340.7844151358.03524413721.020.0144.56											
60         200         120         20         1371.901959         1340.784415         1358.035244         1372         1.02         0.01         44.56											
Means: 3.28 0.07 35.40	60	200	120	20	1371.901959	1340.784415	1358.035244				
								Means:	3.28	0.07	35.40

**Table 5** Objective function data for the 60 instances. All instances have n = 2 and k = 20. The number of complementarities is m. The rank of M and the density of each matrix are indicated. Three relative gaps are given as percentages: (i) the gap between the upper and lower bounds obtained through preprocessing, (ii) the gap between the upper bound obtained from feasibility recovery and the optimal value of the LPCC, and (iii) the improvement in the gap between upper and lower bound effected by the improvement in the LP relaxation obtained through preprocessing.

				Preprocess	Our alg	orithm	default (	PLEX indic	ator constraint
#	m	rank	dense	time	time	nodes	time	nodes	% gap
1	100	30	70	5.93	1.19	404	7.04	2704	/* 8*F
2	100	30	70	4.71	5.70	4340	DNF	6572052	0.485
3	100	30	70	9.35	0.61	260	4.98	1775	0.100
4	100	30	70	1.79	0.13	38	1.70	185	
5	100	30	70	4.80	0.46	210	4.18	1330	
6	100	30	20	1.02	0.02	18	0.73	43	
7	100	30	20	4.52	0.24	44	0.81	103	
8	100	30	20	5.55	0.55	224	5.24	3305	
9	100	30	20	3.56	0.08	32	1.04	204	
10	100	30	20	0.81	0.10	30	1.16	208	
11	100	60	70	1.30	0.24	88	7.08	3940	
12	100	60	70	6.38	0.73	588	384.44	353435	
13	100	60	70	1.21	0.10	20	1.09	49	
14	100	60	70	1.21	1.06	534	9.32	4671	
15	100	60	70	7.25	3.69	1322	DNF	9693517	0.125
16	100	60	20	5.29	1.01	444	2.91	907	
17	100	60	20	6.02	2.06	1324	7.47	5723	
18	100	60	20	1.21	0.35	206	5.55	2672	
19	100	60	20	4.76	0.37	194	3.73	1011	
20	100	60	20	1.00	0.37	226	2.83	708	
			Means:	3.88	0.95	527			
21	150	30	70	24.66	2.40	448	15.58	3307	
22	150	30	70	23.08	0.96	124	4.27	0	
23	150	30	70	5.11	0.16	56	4.58	211	
24	150	30	70	24.49	0.61	98	8.52	2163	
25	150	30	70	24.69	6.17	942	685.24	207429	
26	150	30	20	3.68	0.20	92	3.27	232	
27	150	30	20	16.81	0.20	32	2.76	124	
28	150	30	20	18.82	0.36	78	3.62	370	
29	150	30	20	14.23	1.60	188	7.03	1256	
30	150	30	20	18.49	3.70	682	22.30	10877	
31	150	100	70	26.75	101.60	28538	DNF	13175000	0.146
32	150	100	70	4.86	0.54	192	6.31	580	
33	150	100	70	27.11	29.50	8744	DNF	2323005	0.114
34	150	100	70	14.16	132.03	32124	DNF	1674151	0.206
35	150	100	70	5.21	10.90	2888	DNF	2443005	0.272
36	150	100	20	4.42	16.12	4602	DNF	15955452	0.437
37	150	100	20	22.16	15.45	3064	2338.14	845218	
38	150	100	20	4.84	2.32	584	11.66	1427	
39	150	100	20	22.38	4.84	976	19.74	4403	
40	150	100	20	23.75	32.48	10376	2063.57	1202357	
_			Means:	15.49	16.53	4249			
41	200	30	70	126.63	1546.32	181008	DNF	1368269	0.612
42	200	30	70	12.97	0.22	38	5.50	0	
43	200	30	70	76.18	1.19	66	8.43	189	
44	200	30	70	15.66	1.48	262	29.50	3328	
45	200	30	70	14.40	0.33	64	5.77	0	
46	200	30	20	44.60	0.85	42	5.51	0	
47	200	30	20	49.62	1.17	120	15.08	1714	
48	200	30	20	58.34	22.07	930	1538.63	405689	
49	200	30	20	39.58	0.69	48	5.75	0	
50	200	30	20	48.30	2.80	196	19.23	2749	
51	200	120	70	13.90	176.16	27046	DNF	748850	0.292
52	200	120	70	14.87	319.53	53786	DNF	635814	0.048
53	200	120	70	15.17	25.74	5014	DNF	1721873	0.036
54	200	120	70	13.97	47.07	7736	DNF	1350106	0.028
55	200	120	70	87.88	86.08	8646	DNF	4416595	0.168
56	200	120	20	14.61	283.36	42010	DNF	1371796	0.248
57	200	120	20	82.14	1017.50	97688	DNF	651199	0.736
58	200	120	20	12.90	11.19	1834	DNF	1233196	0.160
59	200	120	20	14.43	31.69	5188	DNF	856913	0.101
60	200	120	20	13.87	491.75	85978	DNF	710977	0.278
			Means:	58.50	203.36	25885			

**Table 6** Performance data for the 60 instances. The final gap obtained by default CPLEX is indicated for the 18 instancesit didn't solve (DNF) within the time limit of 3600 seconds.

	Dimensions		ons		LP	Preprocess	Optimal	% Gap
#	v	b	5115 g	$\operatorname{rank}(Q)$	relaxation	lower bound	value	shrunk
1	50	25	50	25	0.644247	0.67043	0.708016	41.06
2	50	25	50	25	0.691488	0.742312	0.817536	40.32
3	50	25	100	25	0.758156	0.853512	0.90403	65.37
4	50	25	100	25	0.508003	0.719332	1.030281	40.46
5	50	25	150	25	0.634865	0.768284	0.930512	45.13
6	50	25	150	25	0.746323	0.939574	1.179523	44.61
7 8	$50 \\ 50$	$\frac{25}{25}$	$200 \\ 200$	25 25	0.834849	0.882683 0.742780	0.983496	$32.18 \\ 33.90$
9	50	20 50	200	25	0.552456 0.74926	0.742789 0.833068	1.113869 0.974266	37.25
10	50	50	50	25	0.596532	0.680096	0.374200 0.777676	46.13
11	50	50	100	25	0.74352	0.85814	1.025396	40.66
12	50	50	100	25	0.713623	0.897497	1.007106	62.65
13	50	50	150	25	0.734739	0.788571	0.884411	35.97
14	50	50	150	25	0.521193	0.682435	0.990508	34.36
15	50	50	200	25	0.754768	0.782534	0.783561	96.43
16	50	50	200	25	0.754747	0.847708	1.054281	31.04
17	50 50	75 75	50	25	0.649778	0.773429	0.945024	41.88
18	50 50	75 75	50 100	25 25	0.607476 0.720760	0.880724	1.046959	62.17
19 20	$\frac{50}{50}$	$\frac{75}{75}$	$100 \\ 100$	25 25	$0.720769 \\ 0.529814$	$0.812158 \\ 0.667117$	$0.971363 \\ 0.949297$	$36.47 \\ 32.73$
20 21	50 50	75	$100 \\ 150$	25 25	0.529814 0.909594	0.93072	0.949297 0.933821	32.73 87.20
22	50	75	150	25	0.710307	0.836857	1.103089	32.22
23	50	75	200	25	0.718915	0.979733	1.326493	42.93
24	50	75	200	25	0.794803	0.916565	0.950861	78.02
25	50	100	50	25	0.766494	0.894661	1.086423	40.06
26	50	100	50	25	0.485909	0.627154	0.962494	29.64
27	50	100	100	25	0.767284	0.838957	0.95196	38.81
28	50	100	150	25	0.578038	0.699594	0.793066	56.53
29 30	$\frac{50}{50}$	$100 \\ 100$	$150 \\ 200$	25 25	$0.713984 \\ 0.616827$	$0.743964 \\ 0.867557$	$0.760946 \\ 1.064487$	$63.84 \\ 56.01$
31	50	100	200	25	0.651844	0.80751472	0.793559	70.30
32	50	25	50	50	0.644746	0.770222	0.897356	49.67
33	50	25	50	50	0.602	0.759655	0.928742	48.25
34	50	25	100	50	0.660691	0.816037	1.03487	41.52
35	50	25	100	50	0.578159	0.804343	1.031041	49.94
36	50	25	150	50	0.69423	0.856205	1.065649	43.61
37	50	25	150	50	0.790053	0.903077	1.01511	50.22
38 39	$\frac{50}{50}$	$\frac{25}{25}$	$200 \\ 200$	50 50	$0.619887 \\ 0.661197$	$0.766884 \\ 0.876876$	$0.977958 \\ 1.104905$	$41.05 \\ 48.61$
40	50	20 50	200	50	0.57275	0.793187	1.104905	48.01 44.21
41	50	50	50	50	0.54549	0.696252	0.886063	44.27
42	50	50	100	50	0.656456	0.837083	1.073918	43.27
43	50	50	100	50	0.75742	0.844813	0.963771	42.35
44	50	50	150	50	0.67609	0.907669	1.180604	45.90
45	50	50	150	50	0.671718	0.897607	1.117394	50.68
46	50	50	200	50	0.664089	0.816423	1.043968	40.10
47	50	50	200	50	0.571399	0.856775	1.074	56.78
48 49	$\frac{50}{50}$	$\frac{75}{75}$	$\frac{50}{50}$	50 50	$0.514044 \\ 0.647895$	$0.699661 \\ 0.741642$	$0.899591 \\ 0.978757$	$     48.14 \\     28.33 $
49 50	50 50	75 75	100	50 50	0.647895 0.623007	0.741642 0.861603	0.978757 1.058687	28.33 54.76
51	50	75	100	50	0.607434	0.803321	0.940491	58.81
52	50	75	150	50	0.742589	1.022707	1.157709	67.48
53	50	75	150	50	0.706354	0.844787	0.980656	50.47
54	50	75	200	50	0.719345	0.877183	1.064428	45.74
55	50	75	200	50	0.690228	0.845069	0.968196	55.70
56	50	100	50	50	0.642142	0.838403	1.021128	51.79
57	50	100	50	50	0.530841	0.864219	1.13489	55.19
$\frac{58}{59}$	50 50	100	100	50 50	0.647984	$0.832686 \\ 0.925214$	1.119443	39.18
59 60	$50 \\ 50$	$100 \\ 100$	$100 \\ 150$	50 50	$0.70828 \\ 0.544611$	0.925214 0.755445	$1.182871 \\ 1.086004$	45.71 38.94
61	50	100	$150 \\ 150$	50 50	0.71695	0.735445 0.891654	0.996318	62.54
62	50	100	200	50	0.48787	0.771204	1.251992	37.08
63	50	100	200	50	0.64742	0.812777	0.970732	51.14

**Table 7** Values of bilevel instances with dimension of v equal to 50.

	Dir	nensi	ons		LP	Preprocess	Optimal	% Gap	Final	% gaps
#	v	b	g	$\operatorname{rank}(Q)$	relaxation	lower bound	value	shrunk	CPLEX	Our code
64	75	25	50	50	0.70269	0.796622	1.015407	30.04	5.42	
65	75	25	50	50	0.523934	0.662829	0.925725	34.57	solved	
66	75	25	100	50	0.590782	0.727175	1.04984	29.71	17.98	
67	75	25	100	50	0.619884	0.743261	1.056703	28.24	18.16	
68	75	50	50	50	0.561663	0.78647	1.011143	50.01	solved	
69	75	50	50	50	0.560373	0.695674	0.980867	32.18	19.20	
70	75	50	100	50	0.801978	0.889381	1.041901	36.43	solved	
71	75	50	100	50	0.754207	0.881563	0.959749	61.96	solved	
72	75	75	50	50	0.700723	0.815533	1.026476	35.24	7.71	
73	75	75	50	50	0.601257	0.699462	0.930676	29.81	6.95	
74	75	75	100	50	0.346691	0.519392	0.86743	33.16	29.05	
75			50	0.594098	0.738111	1.094845	28.76	25.26		
76	76 100 25 50		50	50	0.882787	0.94306	0.992266	55.05	solved	
77	100	25	50	50	0.567423	0.613049	0.732117	27.70	3.08	
78	100	25	75	50	0.603912	0.639672	0.842811	14.97	20.18	
79	75	25	50	75	0.475835	0.699877			no UB	5.19
80	75	25	50	75	0.524895	0.732925	0.98902	44.82	7.86	
81	75	25	100	75	0.603074	0.783103			16.26	12.05
82	75	25	100	75	0.512566	0.680474			16.44	12.09
83	75	50	50	75	0.546256	0.697274	1.092166	27.66	25.04	
84	75	50	50	75	0.519369	0.68752			18.61	10.53
85	75	50	100	75	0.508331	0.6817			22.22	5.29
86	75	50	100	75	0.619981	0.80781			16.37	4.28
87	75	75	50	75	0.485787	0.630083			27.69	13.95
88	75	75	50	75	0.553843	0.737812			7.31	7.04
89	75	75	100	75	0.416464	0.615691			18.33	13.76
90	75	75	100	75	0.63856	0.783523	1.02771	37.25	solved	

Table 8 Values of bilevel instances with larger dimensions of v. The final gaps obtained by each code are indicated for the instances it did not solve.

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$\begin{array}{c c c c c c c c c c c c c c c c c c c $	#				$\operatorname{rank}(Q)$	Preprocess time				
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$\begin{array}{c c c c c c c c c c c c c c c c c c c $	8	50	25	200	25	25.27	116.34	16582	444.74	245964
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$\begin{array}{c c c c c c c c c c c c c c c c c c c $	16	50	50	200	-					
$\begin{array}{c c c c c c c c c c c c c c c c c c c $	17	50	75	50						
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$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	·				Means:	10.57	15.68	2904		
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$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	39	50	25	200					$\geq 3600$	3371704
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$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	48	50	75	50					43.88	23923
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	62		100	200	50	21.63		323810	$\geq 3600$	
Means: 12.54 472.97 58526	63	50	100	200					353.94	122675
					Means:	12.54	472.97	58526		

**Table 9** Performance on bilevel instances with dimension of v equal to 50.

	Dii	mensi	ons		Preprocess	Our	code	default	CPLEX
#	v	b	g	$\operatorname{rank}(Q)$	time	time	nodes	time	nodes
64	75	25	50	50	11.47	290.70	75616	$\geq 3600$	12455384
65	75	25	50	50	7.73	299.37	68770	$\geq 3600$	17447435
66	75	25	100	50	20.20	396.67	86638	$\geq 3600$	14159416
67	75	25	100	50	16.82	934.03	159358	$\geq 3600$	7263261
				Means:	14.06	480.19	97596		
68	75	50	50	50	12.75	89.09	19502	$\geq 3600$	31475242
69	75	50	50	50	10.46	139.84	28866	$\geq 3600$	1747600
70	75	50	100	50	13.46	51.33	7478	$\geq 3600$	4579441
71	75	50	100	50	17.91	5.75	362	3585.51	1415201
				Means:	13.65	71.50	14052		
72	75	75	50	50	9.32	94.15	16830	$\geq 3600$	4006474
73	75	75	50	50	11.50	86.90	15076	$\geq 3600$	3636937
74	75	75	100	50	21.87	2519.64	267744	$\geq 3600$	1185965
75	75	75	100	50	24.30	1569.69	197996	$\geq 3600$	1144374
				Means:	16.75	1067.60	124412		
76	100	25	50	50	8.99	6.78	408	155.29	27742
77	100	25	50	50	8.18	79.48	5450	$\geq 3600$	2647874
78	100	25	75	50	26.23	1030.75	100044	$\geq 3600$	1049577
				Means:	14.47	372.34	35301		
79	75	25	50	75	8.50	$\geq 3600$	542572	$\geq 3600$	3248691
80	75	25	50	75	11.67	1197.30	212828	$\geq 3600$	2730181
81	75	25	100	75	16.93	$\geq 3600$	449218	$\geq 3600$	13374465
82	75	25	100	75	22.19	$\geq 3600$	268486	$\geq 3600$	13553753
83	75	50	50	75	10.94	3587.62	445828	$\geq 3600$	2126574
84	75	50	50	75	12.60	$\geq 3600$	488384	$\geq 3600$	1758366
85	75	50	100	75	22.05	$\geq 3600$	408186	$\geq 3600$	1269948
86	75	50	100	75	19.24	$\geq 3600$	429360	$\geq 3600$	1389433
87	75	75	50	75	12.25	$\geq 3600$	442694	$\geq 3600$	1984975
88	75	75	50	75	11.15	$\geq 3600$	463366	$\geq 3600$	2227245
89	75	75	100	75	24.20	$\geq 3600$	276876	$\geq 3600$	1390365
90	75	75	100	75	20.69	1221.35	107572	3537.29	1604647

 $\textbf{Table 10} \hspace{0.1 cm} \text{Performance on bilevel instances with larger dimensions of } v.$ 

$\tilde{m}$	$\tilde{n}$	instance	time for our code	CPLEX MIP time	С	PLEX 1	MIP cut	s
					GF	MIR	L&P	IB
100	75	a	303.58	3504.84	5	3	2	
100	75	b	304.58	2686.84	22	1		
100	75	с	60.38	69.09	11			1
100	75	d	328.71	3600.00	19			
100	75	e	94.91	182.61	11	1		
1	mean	or success	218.43	4 of 5				
120	90	a	271.24	1314.80	9	7		
120	90	b	250.26	3600.00	14	2		
120	90	с	126.13	209.95	7	5		
120	90	d	215.09	3600.00	16			
120	90	e	1101.49	3864.48	11	8		
1	mean	or success	392.84	2 of 5				
150	20	a	55.12	702.95	5			
150	20	b	486.00	3263.58	7			
150	20	с	163.54	3319.01	4			
150	20	d	49.13	1260.59	6			
150	20	е	3605.77	3781.06	1			
		success	4 of 5	4 of 5				
200	15	a	154.11	1057.09	7			
200	15	b	81.65	477.30	1			
200	15	с	3604.85	3600.00	4			
200	15	d	128.41	365.59	7		1	
200	15	е	47.71	224.61	5			
		success	4 of 5	4 of 5				
400	5	a	225.54	1573.05	7			
400	5	b	1797.89	3600.00	17			
400	5	с	238.79	183.92	21			
400	5	d	252.45	388.98	20			
400	5	е	47.45	224.18	5			
1	mean	or success	512.42	4 of 5				

Table 11 Performance on 25 inverse quadratic programs. Mean solution time is listed for each set of five problems solved successfully by a code; otherwise, the number of solved instances is given. The number of cutting planes added by CPLEX MIP is also reported; GF are Gomory fractional cuts, MIR are mixed integer rounding cuts, L&P are lift-and-project cuts, and IB are implicit bound cuts.